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NEUTRON FIELD MEASUREMENTS IN PHANTOM WITH FOIL ACTIVATION METHODS

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29 November 1986

Technical Report

CONTRACT No. DNA 001-83-C-0278

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18. SUBJECT TERMS (Continued)

Phantom NBS Standard Source Mid-Thorax Hip

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SECTION 1 INTRODUCTION

The primary goal of this program was to provide a simplified, accurate method of neutron dosimetry for radiobiological studies, so that quantitative assessments of radiobiological effects would not be limited by the accuracy of the dose-measurement method (the cause), but rather by the less reproducible bio-effects (the effect). The probability that such a method will be put to use is a direct function of its accuracy and an inverse function of its costs in turnaround time, manpower, materials, and required expertise or sophistication. The method as presented in Section 4 below incorporates four neutron activation foils with natural thresholds extending from about 0.5 MeV to 8.7 MeV, plus cadmium-covered and bare gold foils $(0.414 \times 10^{-6} \text{ MeV})$ and 0 MeV "thresholds"). It utilizes simple hand-processing of the data and is the order of 6% accurate.

Section 2 presents the "calibration" of the ASTM Standard threshold-foil spectrometry method¹⁻⁴ in the NBS Standard Neutron Field, to establish its accuracy. This method employs the SAND II unfolding code, and was used in AFRRI monkey phantom measurements to provide a basis for evaluating the simple 5-foil (plus cadmium-covered gold) dosimetry method presented in Section 4.

Section 3 presents the results of this same type of spectrum/dose measurement, as carried out at several points inside the monkey phantom. The measurements were carried out both with the standard full dosimetry packet, the "full stack" (i.e., the baseline measurement), and with a much smaller, simplified packet, the "short stack." The accuracy of the short stack, as compared to the full stack, is demonstrated. In this section the dosimetry results, utilizing the SAND II spectrum unfolding code, are compared to the calculated data for absolute spectral flux/dose.

Section 4 presents the development of the simplified "short stack" method and assesses the accuracy of this straightforward, hand-calculational method of threshold-foil neutron dosimetry. The prescription for carrying out accurate dosimetry measurements with the simplified "short stack" method is also presented here.

SECTION 2 CROSS-CALIBRATION IN NBS STANDARD NEUTRON FIELD

A set of seven (7) neutron activation foils was irradiated in the NBS Standard Neutron Field in order to cross-calibrate the threshold foil method of neutron spectrometry. The foils were counted and the data were input to the SAND II unfolding code to obtain the neutron spectrum. A trial spectrum was also input to SAND II. This was the evaluated U-235 fission neutron spectrum discussed below (Figure 1).

Table 1 presents the neutron foil activations, as well as corrections in the NBS Standard Neutron Field for the flux activations gradient and for self-scattering by each foil.

Figure 1 shows the SAND II output spectrum as well as the evaluated U-235 fission neutron spectrum (Designation: XU5-5N1), corresponding to a fluence of 7.786 x 10^{13} n/cm², \pm 1.7%. The corresponding measured value utilizing the ASTM Standard methodology¹⁻⁴ yielded a fluence of 7.98 x 10^{13} n/cm², which is within 2.5% of the NBS value: this is the first and only time the ASTM Method used here has been so evaluated.

Thus, the <u>accuracy</u> of the threshold-foil-spectrometry method, as carried out here, appears to be in the vicinity of 2-3% for a fission-type spectrum. For a water-cooled, water-moderated reactor spectrum, a 1/E tail exists below the fission-like spectrum. However, this 1/E shape is pretty much a canonical appendage; i.e., it is known to rather good accuracy. Therefore, the projected accuracy for a water-moderated reactor is probably the order of 3-4%, based on the NBS calibration results.

It should be pointed out that while the SAND II "calibration" produced a value of neutron fluence within 2.5% of the NBS standard, fluences obtained from some of the individual detectors showed variations approaching 10%. The strength of the multiple-foil approach is apparent here in that those foils departing furthest from the average, after folding in each one with the trial spectrum, are weighted the least. The final spectrum was obtained from the evaluated spectrum (used as input to SAND II) with only one very mild iteration. The spectral-shape constraint, as well as the de-emphasis by the

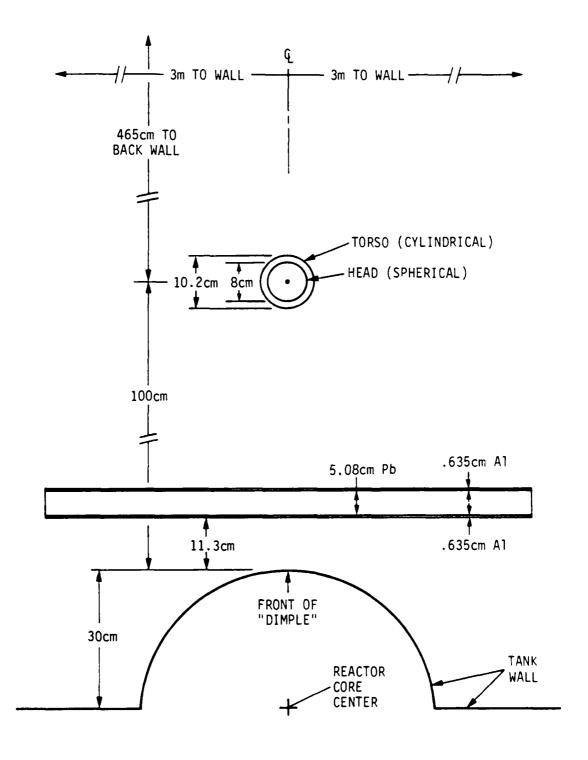


Figure 1. ER1 Exposure Room Top View, Rhesus Monkey Phantom. ER2 is same but without Lead Wall.

Table 1. MBS Cross-Calibration Run Activations Pertinent Parameters & Corrections.

Bol(ce)	1.04-10	6.36-16	2.02-11	1.61-19	8.45-19	1.36-10	6.52-19
,0011100	1.049	1.079	760.0	1.021	1.000	1.049	1.042
,scat	0.972	1.000	0.960	1.008	1.010	1.025	1.021
Roh(en)	1.02-10	5.89-16	11-21.2	1.56-19	8.37-19	1.26-18	6.13-19
-	0.0568	į	0.0588	;	:	:	;
4	0.92	0.50	0.256 0.92 0.96	1.00	0.99	1.00	1.00
Mar-1)	4.126-2	1.546-1	2.256-3 4.126-2 1.723-2	9.532-5	4.051-4	4.621-2	4.621-2
Ey(NeV)	1.596	0.335	0.517 0.743 1.596	0.035	0.010	1.369	1.369
E, th (Nev)	0.50	:	1.45	2.20	2.30	6.3	6.70
101	Np-237	11-115	0-230	Fe-54	11-58	Ng-24	A1-27

No and Noh are given for the fission and non-fission fall activations, respectively

** Read as 4.126 z 10-2

E th . Meutron activation threshold energy

.. . Game rey energy

- Fission Tield

. foil activation (sec-1) uncorrected for scattering and position effects 1. 1(m)

- Fail scattering correction factor; accounts for deviation from neutron-free field due to the presence of the foil 'scat

- foll position correction factor normalized to fluence at the Ni foll (7.79 + 13 ± 1.7%) į

alico) - fall activation (sec-1) corrected for scattering and position effects

code of those foils that are in greatest disagreement in terms of the least-squares fit, result in a multi-foil accuracy considerably greater than that implied by the worst-foil variation. The numerous contributors to the errors in the NBS test include the following.

- 1. Uncertainty in NBS absolute field strength.
- 2. Uncertainty in the flux-gradient and self-absorption corrections.
- 3. Evaluated fission-spectrum-shape uncertainties.
- 4. Errors inherent in histogram-type approximation to the trial spectrum (the evaluated spectrum).
- 5. Threshold-foil cross section uncertainties.
- 6. Gamma-ray-line measurement uncertainties, including NBS-traceable calibration gamma-ray source uncertainties, fluorescent-yield uncertainties, fission branching ratio uncertainties (for fission foils), activation gamma-ray line decay time corrections, and counting statistics.

Efforts were made to reduce many of the uncertainties related to Item 6 above. A large improvement was achieved, as evidenced by the fact that all of the SAND II results, those for the NBS foils as well as those for all the AFRRI phantom spectra presented below, required only one or two mild spectral-change perturbations; a large improvement over earlier measurements of the same kind.

In summary, an accuracy estimate of 3-4% for the AFRRI monkey phantom measurements appears to be quite reasonable. The monkey phantom spectrum differs from the NBS U-235 fission spectrum in that the former has a 1/E tail plus thermal-neutron peak. However, the 1/E tail has a well-known and accurately calculable shape, and only contributes about 11% to the ERI, mid-thorax dose. The thermal peak is not part of this accuracy evaluation, as it can be handled both experimentally and calculationally as an isolated case. (See Section 6 for details.)

SECTION 3 SAND II MEASUREMENTS IN PHANTOM

This section presents a series of six in-phantom spectrum measurements carried out with the short stack of dosimetry foils, one of which was also carried out with the full complement of foils. This will be referred to as the full-stack meeasurement or the basepoint measurement: It is the measurement which serves as a basepoint for the simplified 5-foil dosimetry method presented in Section 4. (See Tables 2 and 3 for the full-stack and the six short-stack foil data, respectively.)

The accuracy of the basepoint measurement was established in Section 2 above where all but the (unnecessary) low-energy neutron-sensitive foils were used for measuring a hard (unmoderated) fission spectrum. In this section, the accuracy of the short stack is evaluated by carrying out the SAND II unfolding both with the short stack and the full stack at the basepoint; i.e., at the mid-thorax point in ER1 (Exposure Room 1), where a 2"-thick lead gamma-ray shield was placed between the reactor and the phantom. (Since 0" to 6" of lead shielding has been used for in-phantom dosimetry measurements, the ER1 exposure turned out to be a good choice for the basepoint measurement. This is especially important for the simplified, hand-calculated dosimetry method developed here and presented in Section 4 below because the 2" lead wall neutron spectrum is intermediate in "softness" between the 0" and 6" lead-shield neutron spectra.)

Figure 2 shows the experimental configuration for the ER1 area, where a 2"-thick lead-wall gamma-ray shield was used. The ER2 configuration was identical, except that no lead shield was used there, and the center of the phantom was 100 cm from the front of the reactor dimple in ER2, instead of 130 cm.

Figure 3 shows a cutaway view, through the rhesus monkey phantom and perpendicular to the reactor-phantom center line, and Figure 4 shows the simple phantom and chair within the plywood enclosure.

Table 2. Basepoint Run (All Foils Present) Activations ER1: Mid-Thorax.

Foi1*	E _n th(MeV)	E ₇ (MeV)	$\frac{\lambda(hr^{-1})}{}$	Yy	Yf	$N_0 \lambda (sec^{-1})^{**}$
197 _{Au}	0+	0.412	1.071-2++	0.95		6.15-16
55 _{Mn}	0+	0.847 1.811	2.687-1 2.687-1	0.99 0.29		3.66-16
235 _U	0+	0.537 0.743 1.596	2.256-3 4.126-2 1.723-2	0.26 0.92 0.96	0.0626 0.0648 0.0626	8.93-11
237 _{Np}	0.50	0.743 1.596	4.236-2 1.723-2	0.92 0.96	0.0568 0.0548	٤.11-12
115 _{In}	1.00	0.335	1.540-1	0.50		2.25-17
238 _U	1.45	0.537 0.743 1.596	2.256-3 4.126-2 1.723-2	0.256 0.92 0.96	0.0588 0.0509 0.0588	7.62-13
232 _{Th}	1.75	0.537 0.743	2.256-3 4.126-2	0.256 0.92	0.0760 0.0534	1.78-13
⁵⁴ Fe	2.20	0.835	9.532-5	1.00	•••	5.87-21
58 _{Ni}	2.90	0.810	4.051-4	0.99	•••	3.20-20
²⁴ Mg	6.30	1.369	4.621-2	1.00	•••	6.08-20
27 _{A1}	8.70	1.369	4.621-2	1.00		3.02-20
90 _{Zr}	14.00	0.910	8.841-3	0.99		2.04-21

^{*} All foils in 0.020" cadmium

 $^{^{\}pm\pm}$ N $_0^{\lambda}$ in sec $^{-1}$ is given for the non-fission foils only N $_0^{o}$ is given for the fission foils

⁺ Since the foils are enclosed in cadmium, the effective threshold for these foils is actually 4.25-7 MeV

⁺⁺ Read as 1.071×10^{-2}

Table 3. Mini-Run (6 Foils Only) Activations Used for Unfolding.

	N _o λ(sec ⁻¹):	ER1	N _o X	(sec ⁻¹): E	R2
Foil*	Thorax **	<u>Head</u>	<u>Hio</u>	<u>Thorax</u>	<u>Head</u>	<u>Hip</u>
197 _{Au} +	6.15-16 ++	5.65-16	6.98-16	1.51-15	1.32-15	1.78-15
115 _{In}	2.25-17	2.10-17	2.86-17	6.59-17	5.57-17	7.6 8-17
54 _{Fe}	5.87-21	5.71-21	7.27-21	1.97-20	1.60-20	2.08-20
58 _{Ni}	3.20-20	3.03-20	3.81-20	1.08-19	8.81-20	1.19-19
24 _{Mg}	6.08-20	5.82-20	6.81-20	2.21-19	1.75-19	2.43-19
27 _{A1}	3.02-20	3.07-20	3.52-20	1.08-19	8.99-20	1.15-19

All foils in .020" cadmium cover

^{**} These activations are taken from the standard (all foils present) run

For a complete list of activation parameters, cf, either Table 1 or 2

⁺⁺ Read as 6.15×10^{-16}

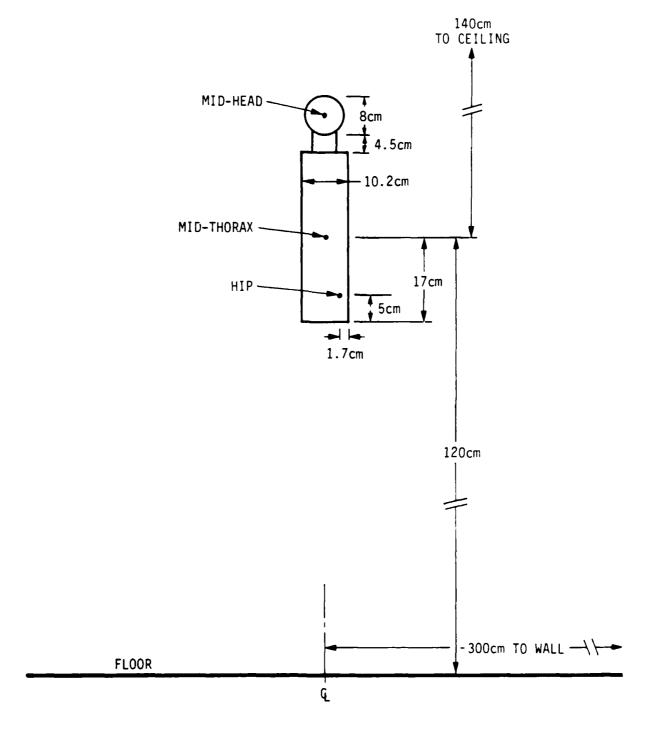


Figure 2. View Perpendicular to Reactor-Phantom Center Line: ER1 and ER2 Exposure Room Configurations.

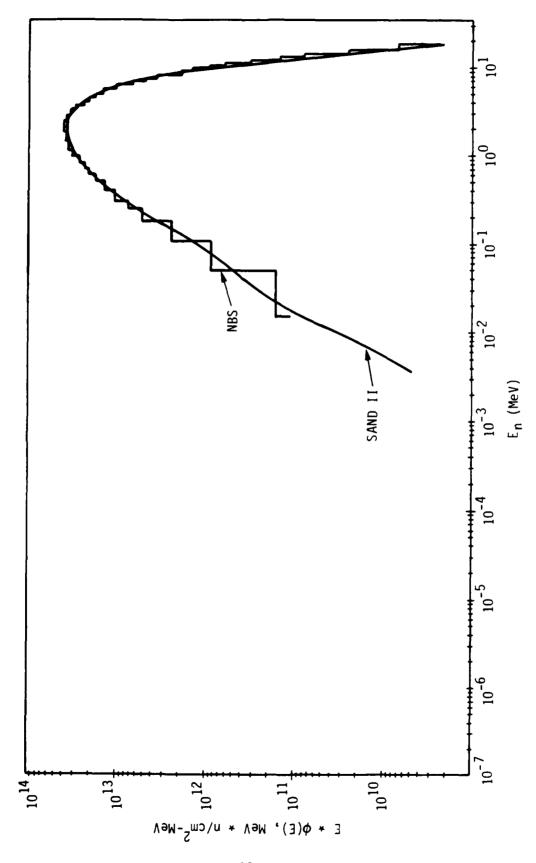


Figure 3. NBS Standard-Neutron-Field Measurement.

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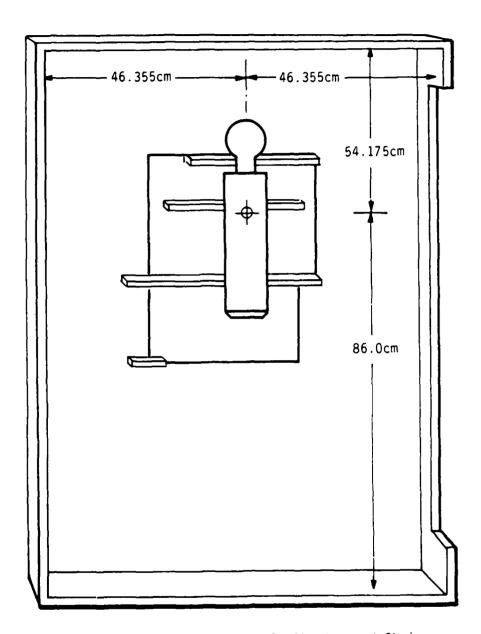


Figure 4. Orientation of Simple Phantom and Chair within the Plywood Enclosure.

The threshold-foil activation data for the full stack of dosimetry foils, the full complement of (ASTM Standard) neutron activation foils, is presented in Table 2 for the rhesus monkey phantom mid-thorax spectrometry/dosimetry exposure in ER1, behind the 2" lead shield; i.e., the basepoint measurement. The threshold-foil data for the six (6) short stack exposures are presented in Table 3, with the ER1 mid-thorax data being common to the Table 2 data. All foils presented in Tables 2 and 3 were cadmium-covered, with a watertight 0.020"-thick cadmium enclosure. The bare gold and manganese foils were exposed in a separate but otherwise identical reactor run, and the activations are not presented in the interest of brevity.

The thermal group is group 37, and the fluence/dose determination is achieved with a simple hand calculation utilizing the gold- or manganese-foil cadmium-difference activations: the methodology of Sections 3 and 4 does not apply to the thermal-neutron fluence/dose. This is adequately covered in Section 6, in terms of the goals of this work.

Table 4 presents a comparison of the SAND II spectrum unfolding code output for the basepoint measurement as carried out with the full stack (standard) and the short stack (labeled "mini"). The calculated spectrum/dose data are also presented in Tables 4, 5 and 6 for calculations carried out in the basepoint location (ER1, mid-thorax), at center head in ER1, and for both mid-thorax and center-head locations in ER2. The comparison of measured and calculated data is presented in Section 3.2 below and includes the experimental normalization of the calculational data, a vital point when considering that one would otherwise have to rely on chamber dosimeters, whose accuracy is, in effect, being evaluated by these more sophisticated spectrometry/dosimetry methods presented here.

The data of Table 4 show an agreement of 2% in total fluence between the full-stack and the short-stack SAND II code outputs. The doses agree to within about 1%. Application of a quicker, analytical technique utilizing these same threshold-foil activation data and the accuracy of that method are presented in Section 4 below. That accuracy assessment is, of course, based directly on the use of the SAND II results (Tables 4, 5 and 6), and on the adequacy of these baseline data.

Table 4. Comparison of Basepoint Run (All Foils Present) with Mini-Run (6 Foils Only) for ER1, Mid-Thorax • Group Fluences (n/cm² • kW • Min), Doses (Rad Tissue).

	Upper	Calc	ulated	Base	point	<u>M</u> ·	ini
Group #	Energy (MeV)	Calc	D _{Calc}	φ _{Meas}	D _{Meas}	φ _{Meas}	D _{Meas}
1	19.6	6.64+3	4.89-5	7.42+3	5.46-5	3.70+3	2.72-5
2	16.9	3.60+4	2.49-4	5.21+4	3.62-4	2.64+4	1.83-4
3	14.9	3.10+4	2.07-4	4.19+4	2.80-4	2.15+4	1.44-4
4	14.2	1.98+4	1.30-4	2.95+4	1.94-4	1.53+4	1.00-4
5 6	13.8	1.42+5	9.02-4	1.73+5	1.10-3	9.35+4	5.95-4
6	12.8	1.30+5	7.95-4	1.55+5	9.49-4	9.66+4	5.91-4
7	12.2	4.99+5	3.03-3	3.52+5	2.14-3	3.47+5	2.11-3
8	11.1	8.55+5	4.90-3	6.45+5	3.70-3	6.43+5	3.69-3
9	10.0	1.51+6	8.35-3	1.10+6	6.06-3	1.09+6	6.01-3
10	9.05	2.37+6	1.25-2	1.79+6	9.43-3	1.78+6	9.37-3
11	8.19	4.37+6	2.24-2	3.23+6	1.65-2	3.20+6	1.64-2
12	7.41	1.23+7	6.08-2	9.05+6	4.48-2	8.99+6	4.45-2
13	6.38	3.32+7	1.53-1	2.69+7	1.24-1	2.74+7	1.26-1
14	4.97	8.02+6	3.51-2	6.89+6	3.02-2	7.07+6	3.09-2
15	4.72	2.86+7	1.23-1	2.35+7	1.01-1	2.41+7	1.04-1
16	4.07	6.75+7	2.75-1	5.87+7	2.39-1	6.07+7	2.47-1
17	3.01	7.74+7	2.66-1	6.39+7	2.20-1	6.64+7	2.28-1
18	2.39	1.34+7	4.22-2	1.10+7	3.47-2	1.14+7	3.60-2
19	2.31	8.19+7	2.54-1	7.03+7	2.19-1	7.38+7	2.29-1
20	1.83	2.02+8	5.41-1	1.61+8	4.31-1	1.66+8	4.46-1
21	1.11	2.30+8	4.70-1	1.90+8	3.90-1	1.90+8	3.89-1
22	.550	1.97+8	2.51-1	1.71+8	2.17-1	1.61+8	2.04-1
23	.158	4.57+7	3.57-2 5.17-2	3.60+7	2.81-2	3.13+7	2.44-2
24 25	.111 .0525	9.51+7 8.11+7	3.17-2 2.49-2	7.16+7 6.04+7	3.90-2 1.85-2	6.31+7 5.32+7	3.43-2 1.63-2
26	.0248	1.05+7	2.18-3	8.23+6	1.72-3	7.22+6	1.51-3
27	.0219	7.93+7	1.14-2	5.60+7	8.01-3	4.96+7	7.11-3
28	.0103	9.04+7	5.52-3	5.00+7 6.93+7	4.23-3	6.22+7	3.80-3
29	3.35-3	8.49+7	1.83-3	6.49+7	1.40-3	5.85+7	1.26-3
30	1.23-3	8.50+7	7.66-4	6.04+7	5.45-4	5.41+7	4.88-4
31	5.83-4	1.72+8	5.36-4	1.24+8	3.87-4	1.10+8	3.43-4
32	1.01-4	1.20+8	1.46-4	8.52+7	1.03-4	7.75+7	9.40-5
33	2.40-5	9.66+7	1.22-4	7.24+7	9.12-5	6.49+7	8.20-5
34	1.07-6	1.51+8	2.96-4	9.91+7	1.94-4	9.12+7	1.78-4
35	3.06-6	8.83+7	2.93-4	6.88+7	2.28-4	6.08+7	2.02-4
36	1.13-6	9.59+7	5.22-4	5.76+7	3.13-4	5.10+7	2.78-4
37	4.14-7	4.34+9	9.72-2	2.84+9	6.37-2	2.84+9	6.37-2
Total		6.60+9	2.76+0	4.57+9	2.25+0	4.48+9	2.28+0

Table 5. ERI Calculated & Measured Group Fluences (n/cm² - kW - Min) and Doses (R/kW - Min) .

			- DIE	d-Thorax			Head	•			
Group .	E _u (NeV)	tele	Mess	DC & 1c	DHeas	Calc	• He a S	Dcalc	Dicess	⁴ ness	Diess
-	19.6	6.64.3	3.70+3	4.89-5	2,72-5	7.97+3	3.89+3	5.87-5	2.86.5	4.2543	11.6
~	•	3.60+4	2.64+4	2.49-4	1,83-4	3.71+4	2.58+4	2.58-4	1.70-4	1.01.6	7 10 7
~	14.9	3.10+4	2.15+4	2.07-4	1.44.1	3.39+4	2.09+4	2.27-4	1.40-4	2.47+4	1.65
-	7.5	1.98.1	1.53+4	1.30-4	1.00-4	1.80+4	1.30+4	1.10-4	1.53-5	1.76+4	1.15-4
·s	13.6	1.42+5	9.35+4	9.05-4	5.95-4	1.35+5	8.02+4	6.61-4	5.11-4	1.07+5	6.12.4
•	12.1	1.30+5	9.66+4	7.95-4	5.91-4	1.12+5	7.86+4	6.83-4	4.81-4	1,11+5	7-67
_	12.2	4.99+5	3.47+5	3.03-3	2.11-3	5.04+5	3.19+5	3.06-3	1.94-3	4.00+5	2.11-1
-	1:1	8.55+5	6.43+5	4.90-3	3.69-3	8.75+5	6.10+5	5.02-3	•	7.40+5	4.25-3
•	0.0	1.51+6	1.09+6	8.35-3	6.01-3	1.59+6	1.08+6	0.75-3	5.96-3	1.26+6	6.94-3
<u>.</u>	9.05	2.37+6	1.78.6	1.25-2	9.37-3	2.90+6	1.95+6	1.53-2	1.03-2	2.04+6	1.08-2
_	- :-	4.37+6	3.20+6	2.24-2	1.64-2	4.93.6	3.34+6	2.52-2	1.71-2	3.70+6	1.89-2
75	7.41	1.23+7	1.99.6	6.08-2	4.45-2	1.27+7	8.71.6	6.28-2	4.31-2	1.05+7	5.20-2
	•	3.32+7	2.74+7	1.53-1	1.26-1	3.14+7	2.55+7	1.45-1	1.18-1	3.32+7	1.53-1
-	1.97	0.02+6	7.07.6	3.51-2	3.09-2	7.76+6	6.59+6	3.40-2	2.00-2	8.61+6	3.77-2
5	•	2.86.7	2.41+7	1.23-1	1.04-1	2.76+7	2.32+7	1.19-1	~	2.93+7	1.26-1
9	4.07	6.75+7	6.07.7	2.75-1	2.47-1	5.79+7	5.32+7	2.36-1	9.60-2	7.39+7	3.01-1
11	0	7.74.7	6.64.7	2.66-1	2.28-1	8.69+7	6.78+7	2.99-1	2.33-1	8.08+7	2.70-1
= :	2.39	1.34+7	1.14.7	4.22-2	3.60-2	1.40+7	1.15+7	4.41-2	3.63-2	1.39+7	4.38-2
•	-	0.19.7	7.38.7	2.54-1	2.29-1	4.06+7	7.48+7	2.92-1	2.32-1	9.02+7	2.81-1
02		2.02+8	1.66+B	5.41-1	1.46-1	1.83+8	1.48+8	4.90-1	3.97-1	2.03+8	5.44-1
12	= :	2.30+8	1.90.8	4.70-1	3.89-1	2.40+8	1.82+8	4.91-1	3.72-1	2.33+8	4.78-1
22	.55	1.97.8	1.61.8	2.51-1	2.04-1	2.30+8	1.88+8	2.92-1	2.39-1	1.91+8	2.43-1
23	. ISB	4.57+7	3.13+7	3.57-2	2.44-2	5.53+7	4.13+7	4.32-2	3.22-2	3.58+7	2.79-2
24		9.51+7	6.31+7	5.11-2	3.43-2	7.86+7	6.35+7	4.27-2	3.45-2	7.16+7	3.90-2
52	.0525	8.11.7	5.32+7	2.49-2	1.63-2	7.61+7	5.77+7	2.33-2	1.77-2	6.04+7	1.85-2
92	.0248	1.05+7	7.22+6	2.18-3	1.51-3	1.25+7	9.74.6	2.61-3	2.03-3	8.22+6	1.72-3
/2	6120	7.93+7	4.96+7	1.14-2	7.11-3	6.53+7	2.75+7	9.37-3	3.94-3	5.65+7	8.08-3
87	.0103	4.044	7+22-9	5.52-3	3.80-3	1.12+8	1.04+B	6.86-3	6.35-3	7.07.7	4.32-3
62	3.35-3	7+67-8	2.85+7	1.03-3	1.26-3	9.53+7	7.30+7	2.06-3	1.57-3	6.63+7	1.43-3
ם י	1.23-3	2.50+7	5.41+7	7.66-4	7-88.7	8.45+7	6.10+7	7.62-4	5.50-4	6.16+7	5.56-4
3	3	1.72+8	3.30+B	5.36-4	3.43-4	1.70+8	1.23+8	5.29-4	3.84-4	1.25+8	3.90-4
35	1.0.1 1.0.1	1.20+B	7.75+7	1.46-4	9.40-5	1.26+8	9.07.7	1.53-4	1.10-4	8.80+7	1.06-4
£ ;	2.40-5	7.66+7	6.49+7	1.22-4	9.50-5	1.00.8	7.31+7	1.27-4	9.24-5	4.69.7	5.91-5
•	6	1.51+8	9.12+7	2.96-4	1.78-4	7.14.8	8.50+7	2.24-4	1.66-4	1.31+8	2.57-4
6	•	8.83+/	/+80.9	2.93-4	2.02-4	9.78+7	7.26+7	3.25-4	2.41-4	6.90.7	2.29-4
4;	1.13-6	9.59+7	5.10+7	5.22-4	2.78-4	1.10•8	6.50+7	5.96-4	3.54-4	5.81+7	3.16-4
`	=	4.34+9	2.84+9	9.72-2	6.37-2	3.32+9	5.96+9	7.44-2	6.62-2	3.70+9	8.28-2

Table 6. ER2 Calculated & Measured Group Fluences (n/cm 2 · kW · Min) and Doses (R/kW · Min) at 100 cm from Reactor Dimple .

			Mid-Thorax	=			Head			*	ē.
Group .	Eu(Nev)	telc.	Meas	DCalc	OHeas	Calc	• He a s	DC+1c	Due . s	*He s	D _{Me e s}
-	19.6	2.23+4	1,30+4	1.64-4	9,57-5	2.70+4	1.13+4	1.99-4	8.32-5	1.41+4	1.04-4
2	16.9	1.24+5	404.6	8.60-4	6.52-4	1.27+5	7.56+4	8.80-4	5.24-4	1,02+5	7.08-4
~	÷.9	1.08+5	7.68.4	7.19-4	5.13-4	1.19+5	6.21+4	7.92-4	4.15-4	8.29+4	5.54-4
•	14.2	6.81.4	5.45+4	4.47-4	3,58-4	6.10•	3.80+4	4.01-4	2.49-4	5.87+4	3.85-4
s	13.8	4.97+5	3.39+5	3.16-3	2.16-3	4.66+5	2.36+5	2.97-3	1.50-3	3.66+5	2,33-3
•	12.8	4.59+5	3.45+5	2.81-3	2.11-3	3.87+5	2.33+5	2.37-3	1.43-3	3.80+5	2.33-3
~	12.2	1.75+6	1.27.6	1.06-2	7.11-3	1.77.6	9.51+5	1.07-2	5.78-3	1.36+6	8.26-3
•	1.1	2.96.6	2.30+6	1.70-2	1.32-2	3.03+6	1.81+6	1.74-2	1.04-2	2.48+6	1.42-2
•	10.0	5.20+6	3.89.6	2.87-2	2.14-2	5.44.6	3.19.6	3.00-2	1.76-2	4.19.6	2.31-2
0	9.05	9.23+6	6.37+6	4.34-2	3.35-2	1.01+7	5.81+6	5.30-2	3.06-2	9+68.9	3.63-2
=	8 .13	1.52+7	1.15+7	7.11-2	5.88-2	1.72+7	1.00.1	8.78-2	5.11-2	1.25+7	6.39-2
12	7.41	4.31+7	3.22+7	2.13-1	1.59-1	4.43+7	2.62+1	2.19-1	1.30-1	3.50+7	1.73-1
13	6.38	1.15.8	9.56+7	5.30-1	4.41-1	1.06+8	7.56+7	4.91-1	3.49-1	1.04+8	4.80-1
=	4.97	2.89+7	2.51+7	1.26-1	1.10-1	2.64+7	1.97+7	1.96-1	1.62-2	2.75+7	1.20-1
7	4.72	1.03+8	6.60+7	1.44-1	3.70-1	9.98+7	6.90+7	4.30-1	2.97-1	9.41+7	4.06-1
•	4.07	2.24+8	2.00.8	9.11-1	8.14-1	1.83+0	1.50+8	7.47-1	6.11-1	2.22+8	9.04-1
-	3.01	2.37.8	2.01+8	8.17-1	6.92-1	2.57+8	1.80+8	8.80-1	6.19-1	2.26+8	7.77-1
•	2.39	3.81+7	3.23+7	1.20-1	1.02-1	4.04+7	2.93+7	1.27-1	9.24-2	3.65+7	1.15-1
-	2.3 1	2.09+B	1.91.8	6.50-1	5.94-1	2.30+8	1.76+8	7.14-1	5.47-1	2.18+8	6.78-1
20	1.63	P. 96.7	4.05+8	1.33+0	1.09.0	4.22.8	3.16+8	1.13+0	8.48-1	4.67+8	1.25+0
21	=:	5.30+B	4.35+8	1.08+0	8.90-1	4.93+8	3.52+8	1.01+0	7.20-1	5.07+B	1.04+0
25	.55	4.17.8	3.59.8	5.30-1	4.56-1	4.46.8	3.77.8	2.66-1	1.79-1	4.22+8	5.36-1
23	158	9.74.7	7.30+7	7.60-2	5.70-2	1.12+8	9.13+7	9.71-2	7.12-2	8.72+7	6.80-2
24	=	2.00.8	7.47.8	1.09-1	7.99-2	1.57+8	8.14.	9.51-2	7.67-2	1.74+8	9.47-2
52	.0525	1.69.6	1.28.8	5.78-2	3.92-2	1.66+8	1.34.8	2-80.5	4.11-2	1.52+8	4.67-2
5	.0248	2.07.7	1.60.7	4.32-3	3.34-3	2.54+7	2.16+7	5.30-3	4.51-3	1.78+7	3.72-3
27	.0219	1.72.8		2.47-2	1.63-2	39+8	8.67.7	1.99-2	1.71-2	1.36+8	1.94-2
28	.0103	1.88+8	1.42+8	1.15-2	8.67-3	2.30+8	1.82+8	1.40-2	1.11-2	1.71+8	1.04-2
5 8	3.35-3	1.82+8	1.38+8	3.92-3	2.98-3	1.94+8	1.63+8	4.19-3	3.51-3	1.61+8	3.48-3
30	1.23-3	1.85+6	1.28+6	1.67-3	7.15-3	8 + 5 9 . 7	1.32.8	1.49-3	1.18-3	1.52+8	1.37-3
]	5.83-4	3.56+8	2.52+8	1.11-3	7.86-4	3.36+8	2.68+B	1.05-3	8.36-4	2.99+8	9.33-4
32	1.0.	2.63.8	1.63+8	3.19-4	7.22-4	2.61+8	2.04+8	3.17-4	2.47-4	2.18+8	2.64-4
33	2.40-5	2.04+8	1.52+8	2.58-4	1.92-4	9 • 96 · 1	1.58+8	2.48-4	2.00-4	1.79+8	2.26-4
7	1.07-5	3.41.6	2.20+8	6.67-4	1.31-4	2.42+B	1.98+8	4.74-4	3.87-4	2.63+8	5.15-4
35	3.06-6	1.90+8	1.46+8	6.32-4	4.85-4	2.10+B	1.71+8	6.97-4	5.68-4	1.71+8	5.68-4
36	1.13-6	2.14+8	1.23+8	1.16-3	6.70-4	2.38+8	1.54+8	1.30-3	6.38-4	1.47+8	8.00-4
~	4.14-7	1.23+10	1.20+10	2.75-1	2.68-1	9.95+9	8.86+9	2.23-1	1.98-1	9.94.9	2.23-1

The measured fluences and doses for the thermal group were determined from the bare and cadmium-covered Au foils activations. SAND II unfolding was not used for this group.

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3.1 SPECTRAL SHAPES.

The SAND II unfolding results of the baseline measurements at the basepoint (ER1, mid-thorax) are shown for the standard dosimetry foil pack in Figure 5 and for the short stack in Figure 6. They are compared to the direct calculational outputs: these calculational data were normalized to some tissue-equivalent (TE) dosimetry measurements taken at some other (distant) time, and the agreement/disagreement with the absolute measurement is therefore misleading. A valid renormalization of the calculated data was done utilizing nickel foils placed on the front of the reactor dimple (see Figure 2), and a discussion of calculational predictions in magnitude is deferred to Section 3.2. The full-stack and short-stack data are in excellent agreement with one another in shape as well as magnitude. The agreement in shape with the calculated (trial) spectrum is also excellent, as evidenced by the fact that the SAND II code carried out its very minimum of only one iteration to obtain the final (measured) spectrum that is in accord with the threshold-foil data.

Figure 7 compares the absolute measured data with the calculated data, as normalized to some TE dosimeter measurement carried out at a different time. The agreement in spectral shape is indeed very good, as it is for the measurement in the rhesus monkey phantom hip in ER1. These data for the hip location, shown in Figure 8, utilized the neutron spectrum calculated at mid-thorax in ER1. The calculated spectrum was arbitrarily fitted to the absolute spectral flux measurement, in Figure 8, to show a good shape comparison.

Figures 9, 10 and 11 show in-phantom spectral data similar to that of Figures 6, 7 and 8, respectively, but for the free-field case (i.e., no lead shield) in ER2. Again, the measured spectral shapes are in good agreement with the calculated shapes.

In Figure 12, a comparison in spectral shape is shown between ER1 and ER2 spectra, mid-thorax. Note that the spectral shapes are practically identical below about 0.4 MeV. This is the "canonical slowing-down spectrum." As normalized in the slowing-down region, the ER1 spectrum above 0.4 MeV is seen

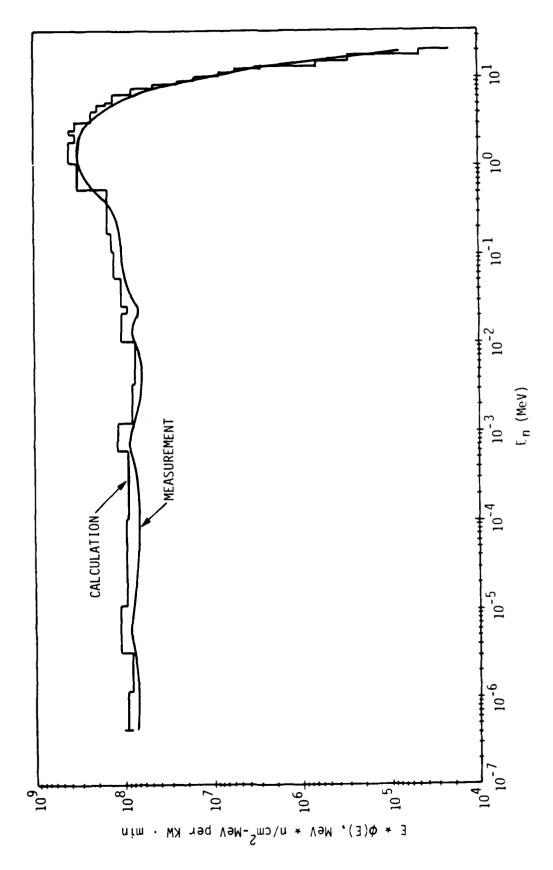
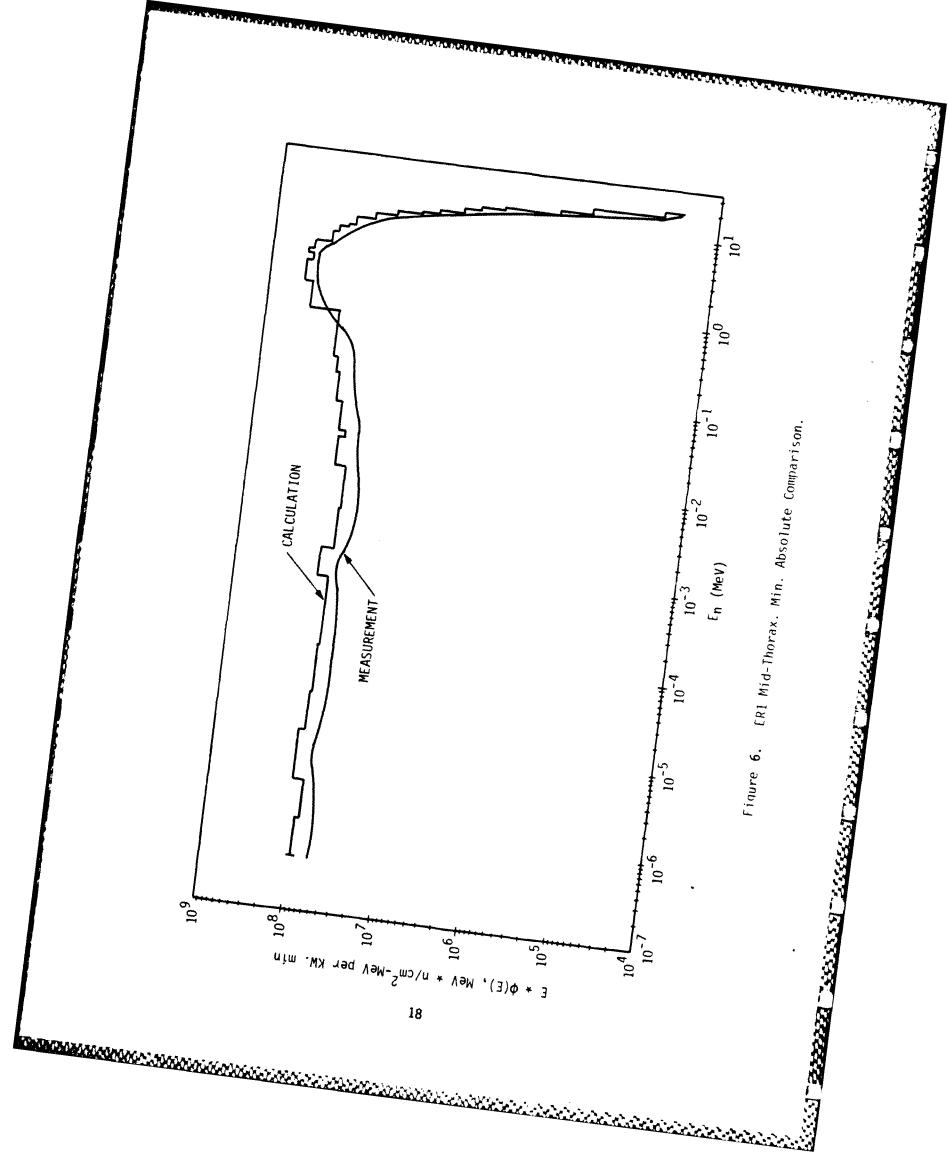


Figure 5. ERI, Mid-Thorax, Max, Absolute Comparison of Calculation with Baseline Measurement.

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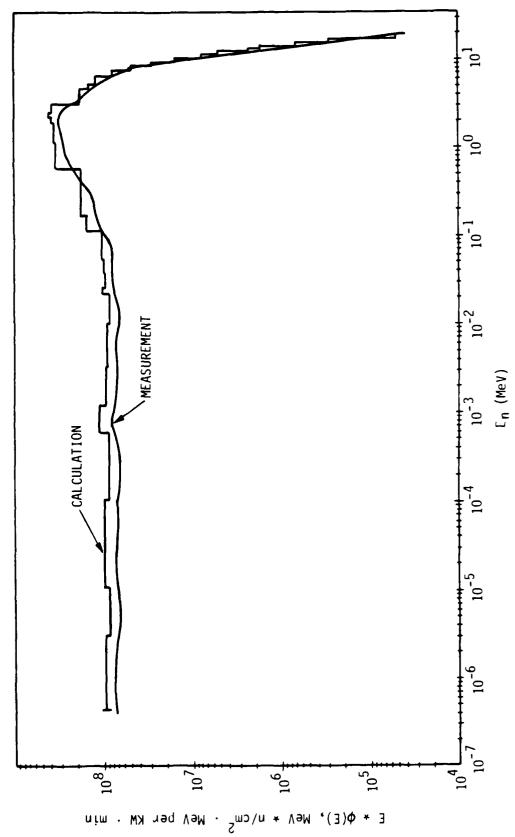


Figure 7. ERl, Head, Absolute Data.

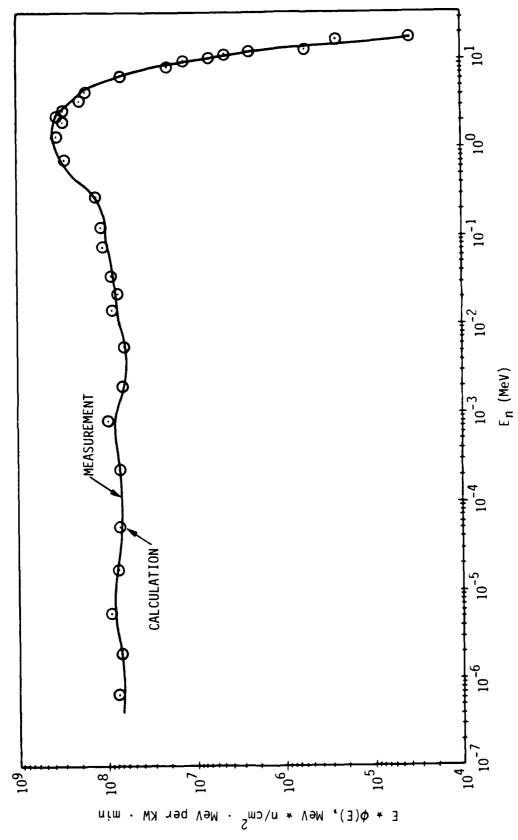


Figure 8. ER1-Hip, Measured, with Calculated Trail-Spectrum (Mid-Thorax) Filtted to it.

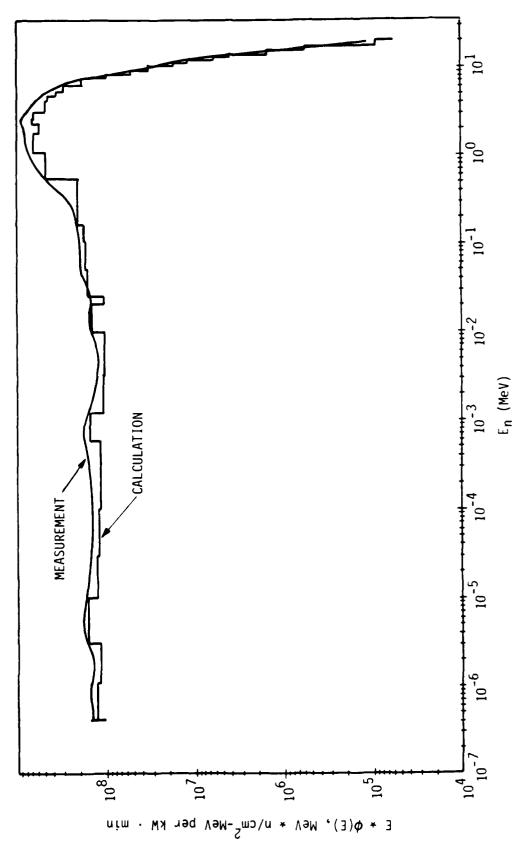


Figure 9. ER2, Mid Thorax, Absolute Fluences.

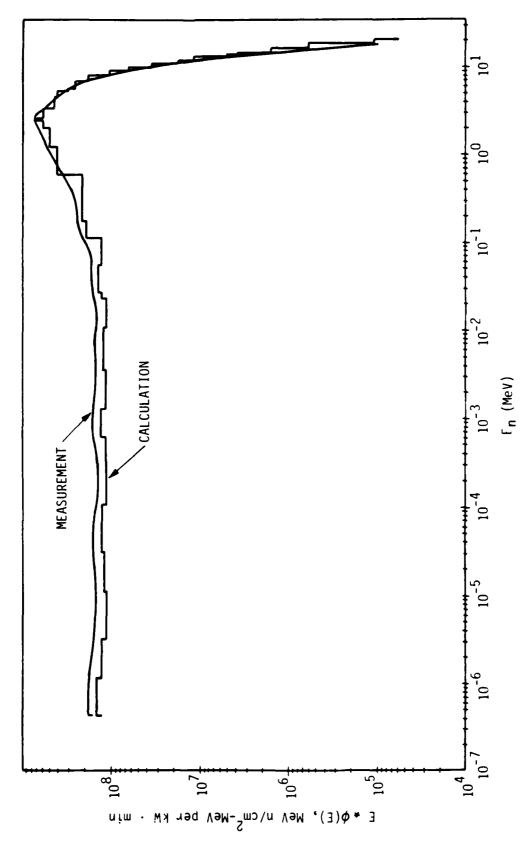
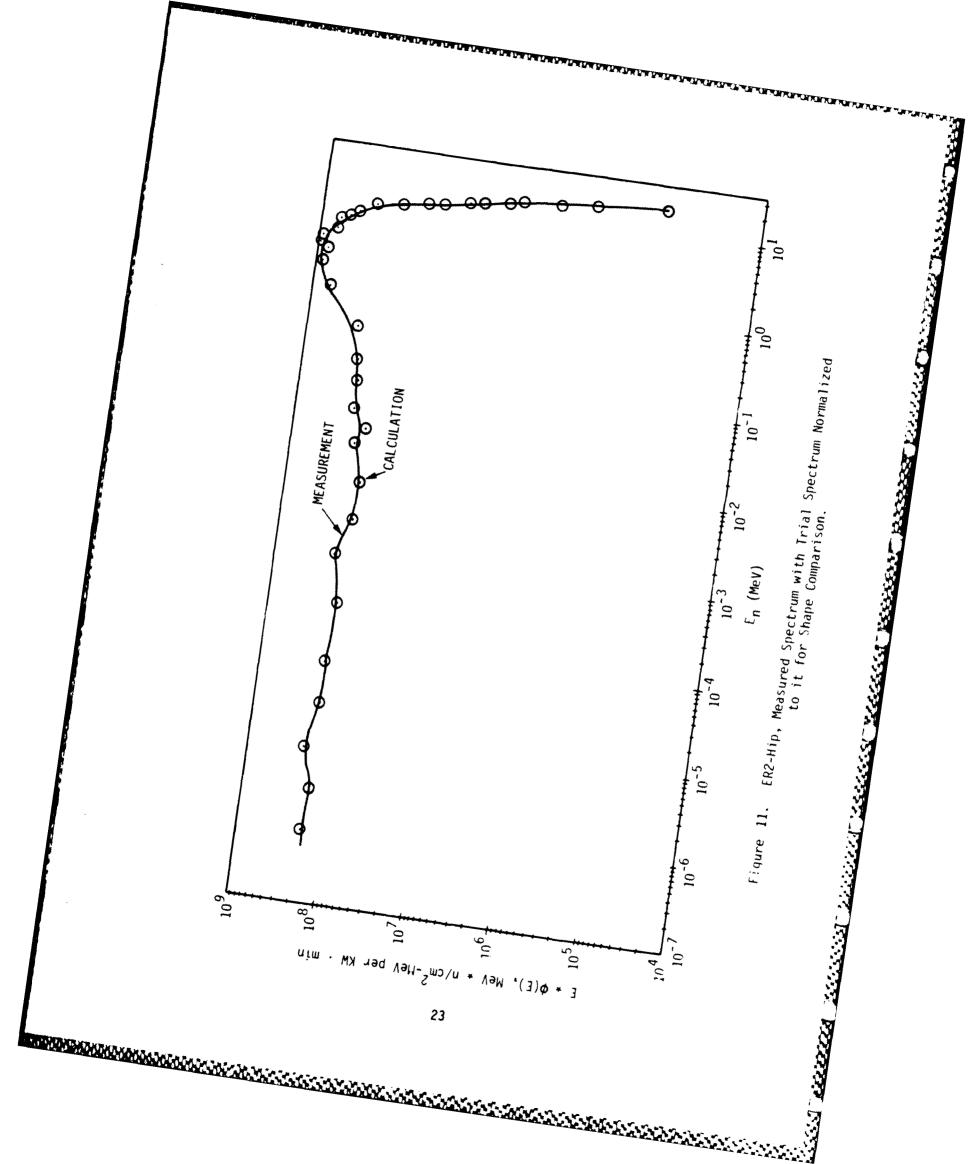
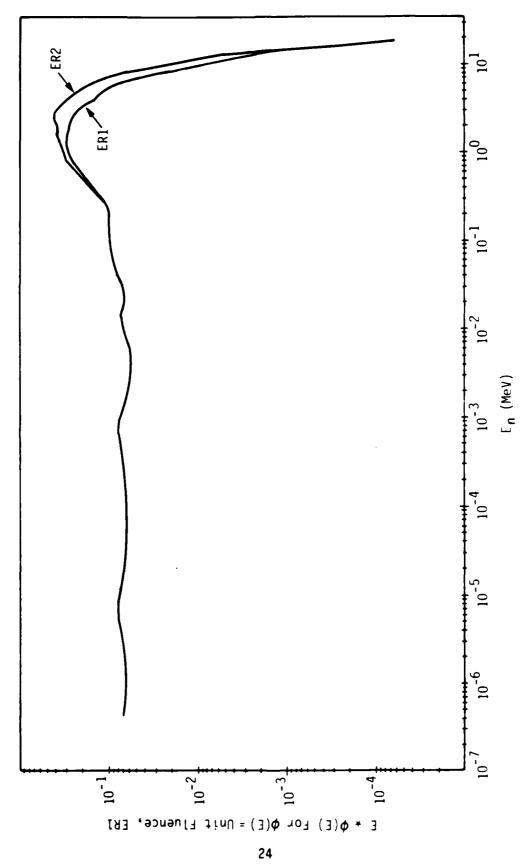


Figure 10. ER2, Head, Absolute Data.

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Spectrum-Shape Comparison, Mid Thorax: Effects of Lead-Wall Shield (ERI). figure 12.

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to be appreciably softer due to the lead wall. Inelastic scattering in the lead shield depletes this high-energy region from 0.4 MeV on up.

3.2 COMPARISONS OF AFRRI ZOOLOGICAL PHANTOM MEASUREMENTS
WITH CALCULATIONS NORMALIZED TO NICKEL FOIL MONITORS.

Neutron fluences have been measured and calculated for mid-head and mid-thorax locations in a simple zoological phantom exposed at the AFRRI TRIGA reactor facility. The calculated data have been produced by Kaul et al, per unit free field KERMA and were first normalized to free field KERMA measured at the intended experiment location, 100 cm from the reactor centerline as follows:

Exposure Room 1 with 2" lead wall 7.8 Rad(tis) per Kw-Min

Exposure Room 2 free field 10.8 Rad(tis) per Kw-Min

The experimental measurements in ER1 were performed at 130 cm from the core centerline rather than 100 cm. Thus, the reported calculations had consequently been adjusted by Verbinski-Cassapakis by dividing by 1.617, a factor determined from measurements made at similar reactors. The calculational results were substantially greater than the measurements, differing by factors ranging from 20 to 100%, which brought to question the practice of normalizing the calculations to standard neutron dosimeters and comparing the results to the much more sophisticated dosimetry measurements presented here.

Because these calculation-measurement discrepancies are much larger than those reported in a similar comparison reported by Kaul et al, it was decided to reexamine the measurement-calculation comparison based on a normalization to nickel foil activation (Ni-58 (n,p) Co-58) data taken at the same time as the phantom measurements. The nickel foils were taped to the aluminum "dimple" which protrudes into each exposure room and into which the reactor is moved at the time of each experiment (see Figure 2). They were positioned at the

reactor midplane along a direct line to the phantom. The nickel foil results, in terms of disintegrations per second per Kw-Min per Ni-58 atom, at exposure time, are as follows:

These values are converted to atoms of Co-58 per Kw-Min per atom of Ni-58 based on a Co-58 disintegration rate of 1.125 ± 10^{-7} per sec., as follows:

The free field fluences on which the phantom calculations are based have been performed using ANISN, 7 a one-dimensional discrete ordinates code, in a spherical geometry model which includes the reactor and all intervening detail. The calculated neutron scalar fluences at the nickel foil locations have been used together with Vitamin-E/ENDF/B-V activation cross sections 8 to calculate the numbers of Co-58 nuclei per Ni-58 nucleus per source neutron which are as follows:

Thus, the numbers of source neutrons per Kw-min required to calculationally reproduce the measured activation in each room are:

Using these values, the neutron KERMA in units of Rad(tis) per Kw-Min at 100 cm from the reactor calculated using one-dimensional methods are:

which are, respectively, 78.6% and 85.8% of the AFRRI monitor values taken at the same locations but on a different day.

The nickel foil-based source normalizations were used to obtain new phantom fluences for comparison with the experimental values. The renormalized calculational results and the experimental results are presented in Table 7. The agreement at the mid-thorax location for neutron energies greater than 10 KeV is 10% or better. Because the mid-head location is actually out of the reactor mid-plane it is expected that the calculated values will show an increase relative to those of the measurements (unless the calculations are corrected for the larger distance of mid-head to reactor center). This is the case for Exposure Room 2 in the bare configuration. However, it is not the case in Room 1, which has the lead shield. The lead appears to have the effect of creating a more uniform exposure over the phantom. For energies below 10 KeV the agreement between calculation and experiment is no better than 10% and is generally much worse. This is most likely the result of using an overly simplified exposure field model.

3.2.1 Summary.

Good agreement has been obtained between calculated and measured fast fluence and KERMA values inside a zoological phantom exposed at the AFRRI TRIGA reactor. This agreement has been obtained on the basis of normalization to nickel activation foil monitor data taken on the face of the reactor containment. Earlier comparisons in which calculations were normalized to paired chamber monitor data taken at the nominal experiment location yielded much poorer agreement. It is expected that better agreement can be obtained at low neutron energies and at locations out of the reactor midplane by using free field fluences obtained by two-dimensional methods, as contrasted to the one-dimensional calculations presented here.

Four-Group Comparison of Calculated Neutron Fluence and Dose with Baseline Measurements Values per Ku-min at 130 cm (ER1) and 100 cm (ER2) from Reactor Centerline. Table 7.

₹	Luci 97		L-PIM	-Thorax	Pi.	Hid-Head	Mid-Thorax	Norax	x Mid-Nead	¥
	(MeV)		N/CB	Rad(tis)	n/cm	Rad(tis)	n/cm ²	Rad(tis)	n/cm ²	Rad(t1s)
1.9	1.96+1	Calc. Neas. c/m	4.09+8 4.53+8 .90	1.38+0 1.53+0 .90	3.95+8 4.26+8 .93	1.34+0 1.44+0 .93	1,31+9 1,29+9 1,02	4.56+0 4.51+0 1.01	1.24+9 1.06+9 1.17	4.34+0 3.70+0 1.17
1.1	1.11+0	Calc. Neas. c/m	5.71+8 5.55+8 1.03	6.54-1 6.77-1	5.78+8 5.70+8 1.01	6.92-1 7.01-1	1.39+9 1.27+9 1.09	1.61+0 1.54+0 1.05	1.32+9	1.56+0 1.41+0 1.11
1.0	1.03-2	Calc. Meas. c/m	7.57+8 6.30+8 1.20	7.60-3 6.73-3 1.13	7.67+8 7.47+8 1.03	8.90-3 9.82-3	1.82+9 1.48+9 1.23	1.82-2 1.56-2 1.17	1.77+9 1.63+9 1.09	2.04-2 1.89-2 1.08
7.E	4.14-7 (Thermal)	Calc. Neas. c/m	3,35+9 2.84+9 1.18	7.49-2 16.37-2 1.18	2.59+9 2.96+9 0.88	5.80-2 6.62-2 0.88	1.05+10 1.20+10 0.88	2.36-1 2.68-1 0.88	8.52+9 8.86+9 0.96	1.91-1 1.97-1 0.96
1		Calc. Meas. c/m	5.09+9 4.48+9 1.14	2.12+0 2.28+0 0.93	4.33+9 4.70+9 0.92	2.10+0 2.22+0 0.95	1,50+10 1,60+10 0.94	6.42+0 6.33+0 1.01	1.29+10 1.28+10 1.01	6.11+0 5.33+0 1.15
) N	*Normalized to		Nickel Foil mo	monitors on containment	ontainnent	face.				

SECTION 4 THE 5-FOIL NEUTRON DOSIMETRY METHOD

This section presents a simple hand-calculational (perturbation) method of threshold-foil dosimetry in which 5 threshold-activation foils are used. (The SAND II code is not required; neither is there need for a trial spectrum. However, the "incremental doses" assigned to each foil (Table 8) were derived from the SAND II basepoint measurement.) One foil, the nickel foil, is used to derive the first-order dosimetry value, $D^{O}(Ni)$. This is simply equal to D^{B} , the dose of the baseline run (i.e., the "standard" run in Table 4), multiplied by the ratio of the nickel-foil activations.

$$D^{O}(Ni) = D^{B} R(Ni)$$
 4.1

where $R(Ni) = (N_o \lambda)^X (Ni)/(N_o \lambda)^B (Ni)$.

Here, $(N_0\lambda)^X$ refers to the dosimetry measurement under consideration and $(N_0\lambda)^B$ refers to the baseline-run value (the value for mid-thorax, ER1, in Table 3).

Now, if there were no change of spectral shape between the baseline run and run x, Equation 4.1 would give an accurate measure of the dose: The nickel foil has proven to be, by far, the most reliable foil in terms of reproducible results.

The nickel foil by itself cannot, of course, correct for a change of spectral shape. However, the foils with thresholds at energies different from the nickel threshold of about 3 MeV do provide change-of-spectral-shape data. This is done by means of the following "shape-perturbation" correction.

$$D^{X} = D^{O}(Ni) + \sum_{i} (A_{i}F_{i}-1)$$
 4.2

Here, the A_i refers to the activation ratios of the i-th foil (i.e., Au, In, Fe and Mg) for the unknown and the baseline run, divided by R(Ni): i.e.,

$$A(Au) = R(Au)/R(Ni)$$

where

$$R(Au) = (N_0 \lambda)^X (Au)/(N_0 \lambda)^B (Au)$$
.

The various $\mathbf{F_i}$ refer to the fraction of the dose "sensed" by the i-th foil, in the case of the baseline run, excluding the nickel foil. The $\mathbf{F_i}$ values were determined from the dose values $\mathbf{D_{STD}}(R)$ given in Table 8 for the "standard" or baseline run; i.e., the ERI, mid-thorax, full-stack spectrum. The corresponding energy range (region of maximum influence) for each foil is also presented in Table 8. Table 9 presents the value of $\mathbf{F_i}$ for each of the 4 "shaping-sensing" foils.

4.1 APPLICATION OF THE SIMPLE 5-FOIL DOSIMETRY METHOD.

The foil-activation values presented in Table 3, for all six (6) spectrum measurements, were utilized to provide the doses in phantom by means of Equations 4.1 and 4.2 above. The dose-fractions F_i , Equation 4.2, are given in Table 9 for each "shape-correction" foil i (i.e., Au, In, Fe and Mg).

The values of A_i (i.e., R(Mg)/R(Ni)) are presented in Table 9, along with $D^0(Ni)$, the first order nickel-foil dose determination, and D^X , the second order determination in which a change-of-spectral-shape correction has been applied. Note how the A_i systematically increase with neutron energy for the ER2 spectra: As shown in Figure 11, these spectra are appreciably harder than the ER1 spectra, due to the down-scattering (in energy) of neutrons by the 2-inch thick lead wall in ER1.

The values of D^X are presented in Table 9, along with D^S , the values derived from the SAND II unfolding-code spectra. The last row presents the percent variation of D^X from D^S . The standard deviation for the ER1 (excluding mid-thorax) and ER2 data are 6%.

Table 8. Standard Run Parameters for Short-Stack Dosimetry at AFRRI.

<u>Foil</u>		sented y Range E _U (MeV)	$(N_o \lambda)_{STD}(sec^{-1})$	N ⁺ STD	D _{STD} (R) ⁺⁺
197 _{Au}	4.14-7	0.55	6.68-16	2.09+4	1460.2
115 _{In}	0.55	2.31	2.25-17	7.03+2	4551.4
54 _{Fe}	2.31	3.01	5.87-21	1.83-1	1114.6
58 _{N i}	3.01	6.38	3.20-20	1	2142.9
24 _{Mg}	6.38	19.6	6.08-20	1.90+0	375.4

^{*} The phantom mid-thorax exposure behind 2" lead at ER1 is taken to be the standard run.

 $N^{+}_{\mbox{\scriptsize STD}}$ is the Ni-normalized foil activation.

 $^{^{++}}$ D_{STD}(R) are doses for the 4187.5 kW \cdot Min exposure, between E_L and E_V. Columns 2 and 3, taken from the basepoint ("standard") measurement, Table 4, but with ICRP rather than ICRU dose conversions. The ICRP vs. ICRU differences are irrelevant here.

Compared to D_s from SAND 11 Unfolding with 5 Folls. Doses in Rad (Tissue) per kN· Min. Table 9. Bose Determinations via 5-foil Bosimetry Stack for Epithermal Newtrons, per kW·Min, The Bo and DA Values are Hand-Calculated Values (See Equations 4.1 and 4.2).

		ER1 (2" Lead)	(pra		=	ER2 (No Lead)	
	<u>.</u> -	Mid-Thorax	Head	d I	Hid-Thorax	700	E C
Ds (Epithermal)	1	2.22	2.15	2.81	90.9	5.13	7.21
R(Au)/R(B1)	0.20	1.00	0.970	0.954	0.746	0.780	0.778
R(1n)/R(H1)	0.60	1.00	0.986	1.067	0.867	0.900	0.917
R(Fe)/R(N1)	0.15	1.00	1.027	1.041	0.993	0.991	0.953
R(Hg)/R(H1)	9.09	1.00	1.011	0.941	1.075	1.047	1.074
R(B1)		1.000	0.98	1.19	3.38	3.75	3.72
B*(#1)		2.22	2.10	2.64	7.49	6.11	8.26
K O		2.22	5.08	2.60	6.53	8.48	7.45
6(S)(8f-85)		:	-3.4	-8.0	1.1	9.9 +	+3.4

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The 5-foil stacks for the mid-thorax spectrum, ER1, were taken from the baseline-run data presented in Table 2. When the SAND II code was run with the 5-foil stack, the resulting epi-thermal neutron dose was about 1% higher. This is somewhat representative of the error resulting from utilizing a small fraction of the full-foil-stack data. This is to be compared with the 6% standard deviation resulting from utilizing the same five foils, but without SAND II.

SECTION 5

COMPARISON OF SAND II (5-FOIL STACK) AND THE SIMPLE 5-FOIL DOSIMETRY METHOD PRESENTED IN SECTION 4

The improvement, due to the use of SAND II, derives from making full use of the a priori knowledge of the spectral shape. This a priori knowledge is input to SAND II in the form of the trial spectrum; i.e., the spectrum from the best available calculation. Thus, SAND II also represents a perturbation method of neutron spectrometry/dosimetry. The best-known (input) spectrum is perturbed in shape and absolutely adjusted in magnitude to agree with the foil-activation data.

The 5-foil dosimetry method presented in Section 4 can be thought of as a perturbation method as well in that it corrects for variations in spectral shape via an approximate 5-group ("histogram") approach. By contrast, SAND II makes full utilization of the a priori knowledge of the spectral shape. However, the estimated 6% accuracy for the 5-foil 5-group "histogram" method may be adequately accurate, compared to the 3-4% estimated accuracy for the maxi-foil dosimetry stack used with SAND II. Consider further that no accurate a priori knowledge of spectral shape is needed for the 5-foil 5-group "histogram" method of Section 4, making it much easier to apply.

The estimated 6% accuracy is to be contrasted to the inaccuracies implied in Section 3.2, where the calculational results were first obtained via normalization to measured free field KERMA (measured with TE dosimeters). Discrepancies ranging from 20 to 100% were reduced to the order of 10%, according to Section 3.2. This section, as well as Section 3.2.1, was taken directly from a communication by Dean Kaul (Letter, August 22, 1985).

SECTION 6 THERMAL (SUB-CADMIUM) NEUTRON DOSE

The thermal-neutron dose is derived from gold foil data, utilizing the cadmium-difference method. The gold-foil cross section is taken as 78.54 barns, which is the cross section derived from a Maxwellian thermal-neutron peak plus 1/E tail extending out to 0.414 eV. This cross section value is more accurate than the value of 98.8 barns, for 2200 m/sec neutrons.

A value of 2.24 \times 10⁻¹¹ rads per unit neutron fluence is the proper dose conversion factor (ICRU muscle KERMA).

It must be pointed out that the bare and cadmium-covered gold (or manganese or cobalt) foils cannot be exposed one behind the other, or even side-by-side, because of the cadmium-produced local flux depression: If side-by-side, but well separated (say by 3 or 4 cadmium "diameters"), then it is incumbent on the experimenter to prove that both are in identical epicadmium neutron-flux fields. In the measurements presented here, the bare foils were exposed in separate reactor runs, with run-to-run normalization carried out with nickel activation foils taped to the reactor entrance window (on the front of the "dimple" in Figure 2).

These nickel monitor foils proved to be a fortunate choice because they provided a basepoint for the "reactor power" or "reactor source term" renormalization. This is an otherwise formidable task, requiring flux plots throughout the reactor, an exact knowledge of day-to-day secondary standard (fission chamber, say) calibration, of the thickness of water between reactor and exit window to ERI/ER2, etc. (Prior to this, the calculation was renormalized to some dosimeters used at a much different time. This proved unfortunate, as can be seen from the results of Table 4, for example. Renormalization to the nickel foils on the dimple provided a meaningful comparison of measurement and calculation, which would have otherwise proven untenable.)

SECTION 7

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